



# wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 6, 2026 – 07:32 PM UTC

PDB ID : 5A2N / pdb\_00005a2n  
Title : Crystal structure of the nitrate transporter NRT1.1 from Arabidopsis thaliana.  
Authors : Parker, J.L.; Newstead, S.  
Deposited on : 2015-05-20  
Resolution : 3.70 Å(reported)

This is a wwPDB X-ray Structure Validation Summary Report for a publicly released PDB entry.

We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0  
Xtrriage (Phenix) : 2.0  
EDS : 3.0  
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)  
CCP4 : 9.0.010 (Gargrove)  
Density-Fitness : 1.0.12  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.49

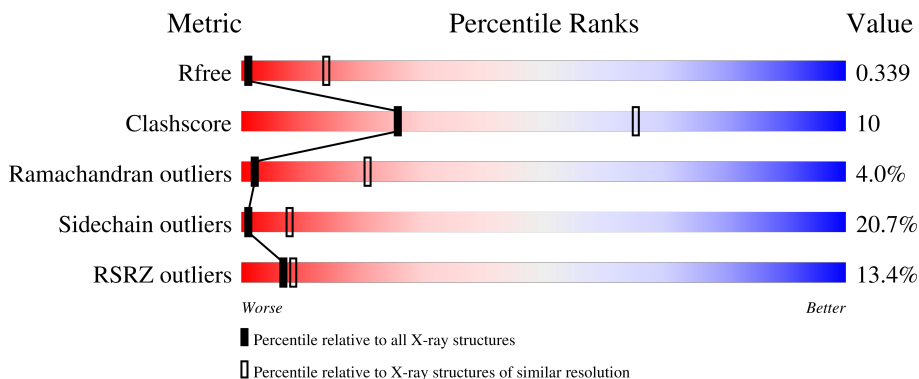
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 3.70 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	180053	1131 (3.80-3.60)
Clashscore	190562	1171 (3.80-3.60)
Ramachandran outliers	187476	1129 (3.80-3.60)
Sidechain outliers	187428	1126 (3.80-3.60)
RSRZ outliers	180081	1130 (3.80-3.60)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	590	
1	B	590	

## 2 Entry composition

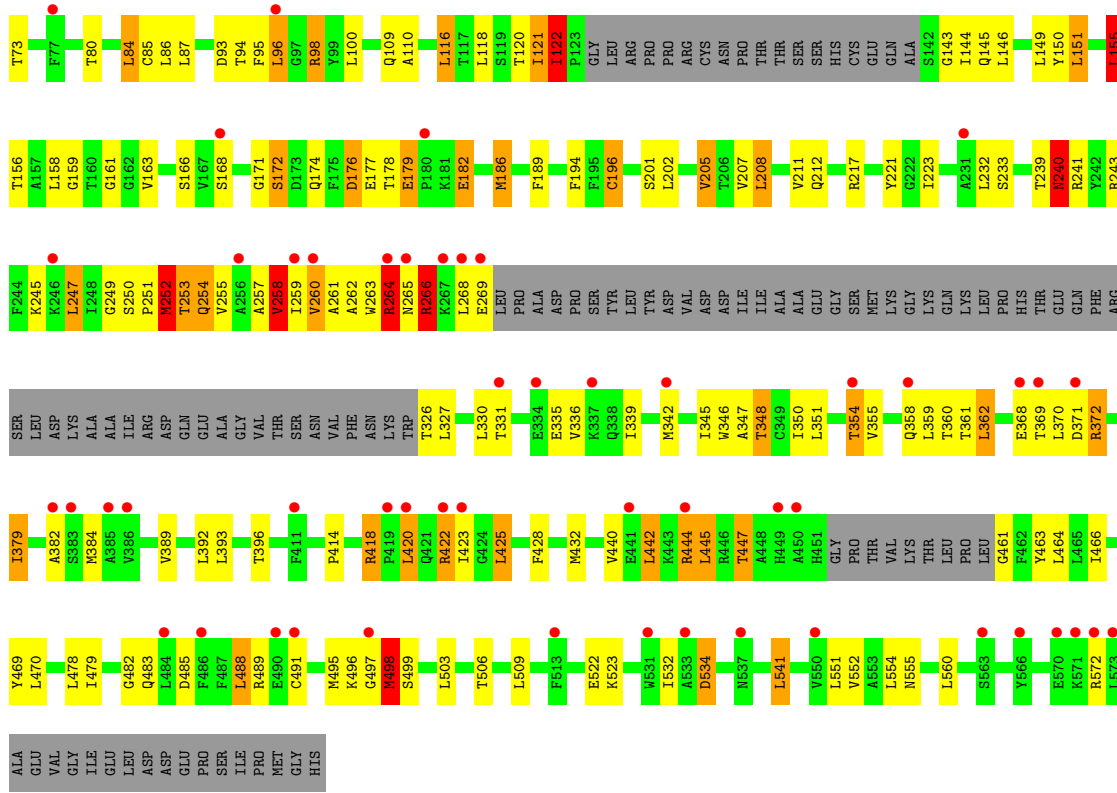
There is only 1 type of molecule in this entry. The entry contains 7474 atoms, of which 166 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called PROTEIN NRT1/ PTR FAMILY 6.3.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
1	A	470	3737	2408	83	599	629	18	0	0	0
1	B	470	3737	2408	83	599	629	18	0	0	0





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	121.80Å 124.03Å 153.32Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	27.53 – 3.70 27.53 – 3.70	Depositor EDS
% Data completeness (in resolution range)	86.3 (27.53-3.70) 86.0 (27.53-3.70)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.34 (at 3.74Å)	Xtrriage
Refinement program	BUSTER 2.10.0	Depositor
R, $R_{free}$	0.278 , 0.310 0.307 , 0.339	Depositor DCC
$R_{free}$ test set	1105 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	105.4	Xtrriage
Anisotropy	0.025	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.24 , 85.9	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.44$ , $\langle L^2 \rangle = 0.26$	Xtrriage
Estimated twinning fraction	0.037 for k,h,-l	Xtrriage
$F_o, F_c$ correlation	0.83	EDS
Total number of atoms	7474	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	81.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.81% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 5$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z  > 5$	RMSZ	$\# Z  > 5$
1	A	0.86	2/3734 (0.1%)	1.57	46/5067 (0.9%)
1	B	0.88	2/3734 (0.1%)	1.61	57/5067 (1.1%)
All	All	0.87	4/7468 (0.1%)	1.59	103/10134 (1.0%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	252	MET	SD-CE	6.63	1.96	1.79
1	A	122	ILE	CA-C	5.53	1.58	1.52
1	A	466	ILE	CA-CB	5.19	1.57	1.53
1	B	266	ARG	CA-C	5.15	1.59	1.52

The worst 5 of 103 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	498	MET	N-CA-C	-9.60	99.40	112.94
1	A	500	THR	N-CA-C	-8.86	102.46	113.18
1	B	24	ASP	CA-CB-CG	7.88	120.48	112.60
1	A	95	PHE	N-CA-C	7.49	120.63	108.34
1	A	24	ASP	CA-CB-CG	7.45	120.05	112.60

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3654	83	3786	68	0
1	B	3654	83	3786	74	0
All	All	7308	166	7572	142	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 10.

The worst 5 of 142 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:263:TRP:HA	1:B:266:ARG:HB3	1.67	0.76
1:B:250:SER:HB2	1:B:251:PRO:HA	1.67	0.75
1:B:251:PRO:O	1:B:254:GLN:HG3	1.87	0.74
1:A:108:ILE:HG21	1:A:158:LEU:HD23	1.68	0.73
1:B:57:THR:HB	1:B:217:ARG:HH22	1.56	0.71

There are no symmetry-related clashes.

## 5.3 Torsion angles [i](#)

### 5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	462/590 (78%)	391 (85%)	48 (10%)	23 (5%)	1	17
1	B	462/590 (78%)	388 (84%)	60 (13%)	14 (3%)	3	26
All	All	924/1180 (78%)	779 (84%)	108 (12%)	37 (4%)	2	21

5 of 37 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	55	LEU
1	A	252	MET
1	A	411	PHE

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Mol	Chain	Res	Type
1	A	413	TYR
1	A	419	PRO

### 5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	386/489 (79%)	311 (81%)	75 (19%)	1	9
1	B	386/489 (79%)	301 (78%)	85 (22%)	1	7
All	All	772/978 (79%)	612 (79%)	160 (21%)	1	8

5 of 160 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	254	GLN
1	B	470	LEU
1	B	264	ARG
1	B	360	THR
1	B	506	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 12 such sidechains are listed below:

Mol	Chain	Res	Type
1	A	537	ASN
1	A	555	ASN
1	B	483	GLN
1	B	67	ASN
1	A	240	ASN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

#### 5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

#### 5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

#### 5.6 Ligand geometry [i](#)

There are no ligands in this entry.

#### 5.7 Other polymers [i](#)

There are no such residues in this entry.

#### 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

## 6 Fit of model and data [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled '#RSRZ > 2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled 'Q < 0.9' lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	A	470/590 (79%)	0.61	66 (14%) <b>6</b> <b>8</b>	3, 66, 230, 270	0
1	B	470/590 (79%)	0.66	60 (12%) <b>7</b> <b>9</b>	3, 57, 204, 239	0
All	All	940/1180 (79%)	0.63	126 (13%) <b>7</b> <b>8</b>	3, 60, 214, 270	0

The worst 5 of 126 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	21	ARG	7.8
1	B	486	PHE	5.8
1	A	329	THR	5.7
1	B	450	ALA	5.2
1	A	96	LEU	5.2

### 6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

### 6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

### 6.4 Ligands [i](#)

There are no ligands in this entry.

### 6.5 Other polymers [i](#)

There are no such residues in this entry.